

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101218\
 Data File : BG037294.D
 Acq On : 12 Oct 2018 17:01
 Operator : JU/SJ
 Sample : PB113784BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113784BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101118.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.198	319	325	335	rVB2	30423	49565	1.31%	0.279%
2	5.656	396	403	414	rBV	775784	1294049	34.30%	7.290%
3	7.254	667	675	684	rBV	824708	1425370	37.78%	8.030%
4	7.642	732	741	749	rBV	754205	1330662	35.27%	7.496%
5	8.118	813	822	831	rVB	162463	285070	7.56%	1.606%
6	8.441	868	877	887	rVB	565504	1007208	26.70%	5.674%
7	9.293	1014	1022	1034	rBV	448356	820520	21.75%	4.622%
8	10.938	1295	1302	1311	rBV	221076	416178	11.03%	2.344%
9	13.377	1710	1717	1727	rVB	1318882	2105693	55.82%	11.862%
10	14.751	1942	1951	1959	rBV2	391730	640085	16.97%	3.606%
11	16.238	2197	2204	2220	rBV	1004224	1601525	42.45%	9.022%
12	17.495	2412	2418	2425	rBV2	576556	918874	24.36%	5.176%
13	20.098	2855	2861	2868	rBV	2750731	3772430	100.00%	21.251%
14	21.784	3141	3148	3158	rBV	624166	1052717	27.91%	5.930%
15	25.133	3707	3718	3736	rVB2	340448	1031474	27.34%	5.811%

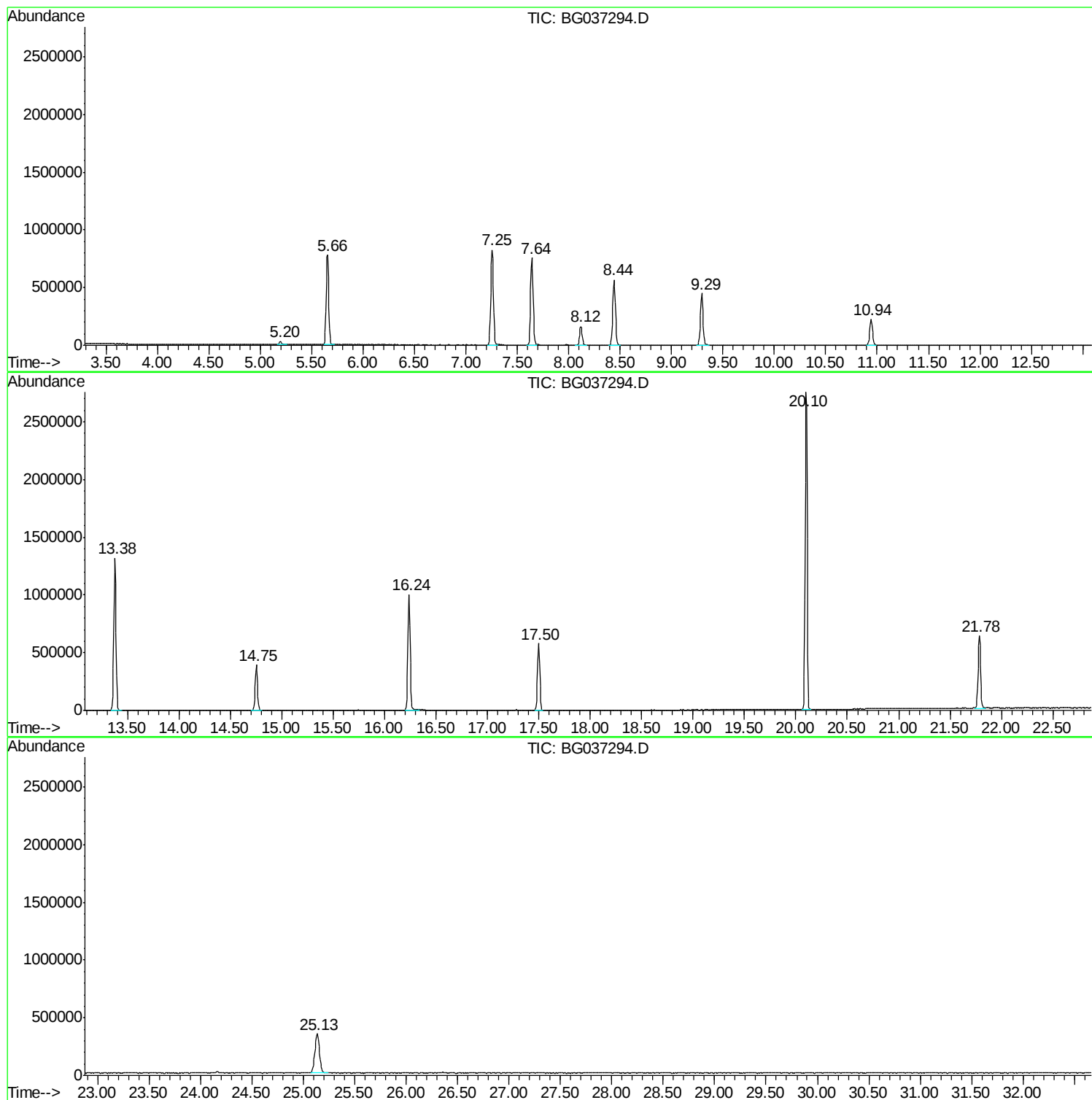
Sum of corrected areas: 17751420

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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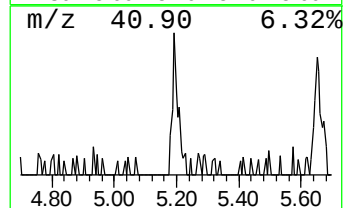
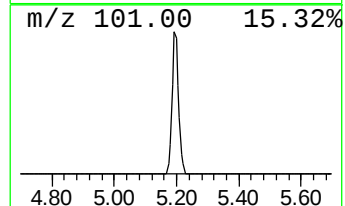
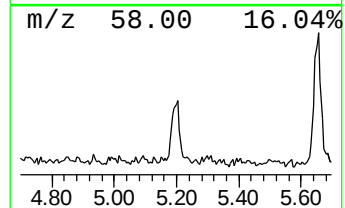
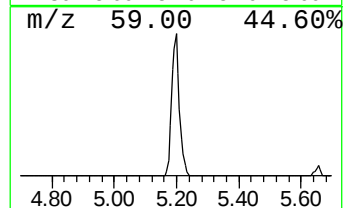
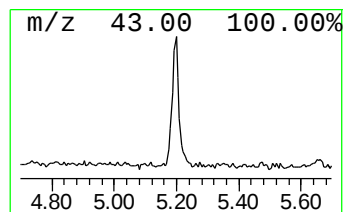
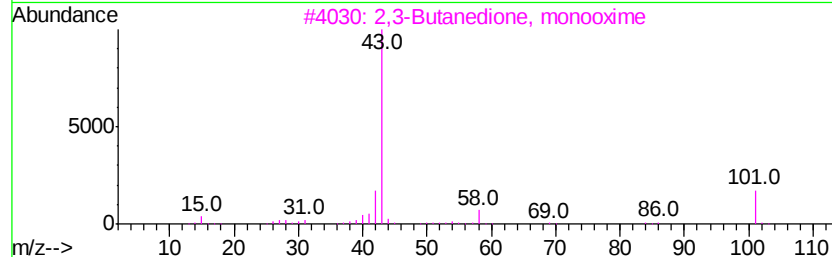
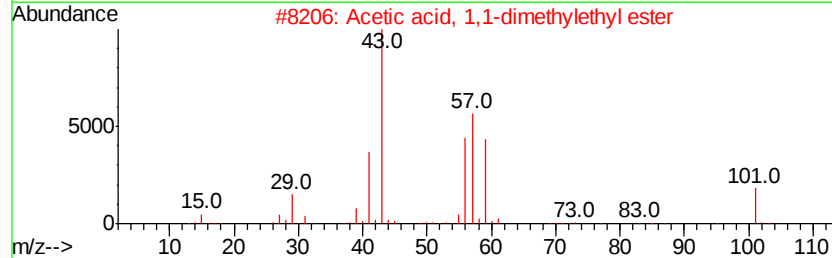
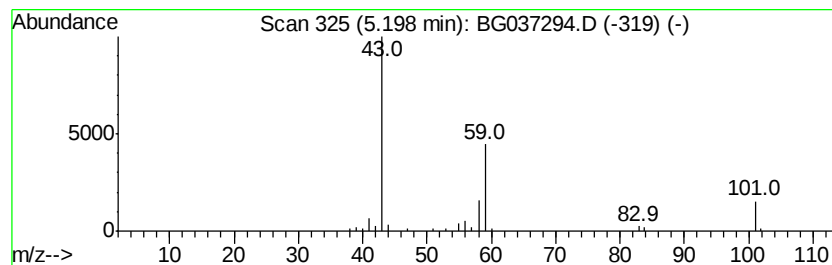
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	3.48 ng	49565	1,4-Dichlorobenzene-d4	8.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27	
4	Ethyl ether	74	C4H10O	000060-29-7	9	
5	Butane	58	C4H10	000106-97-8	7	



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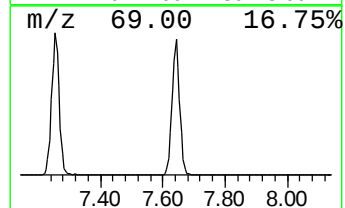
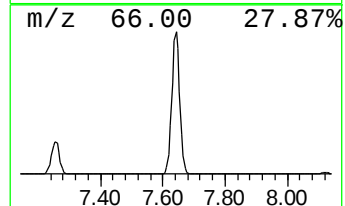
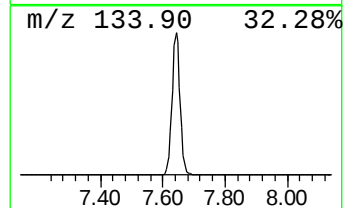
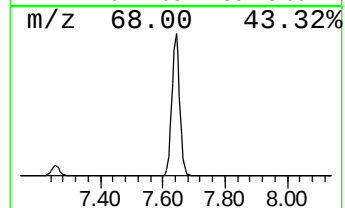
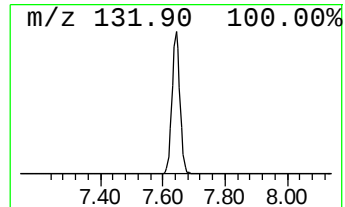
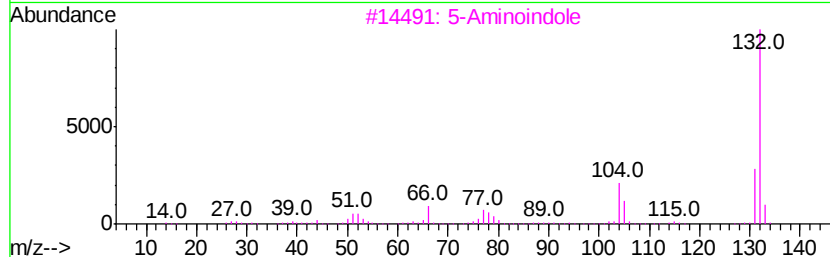
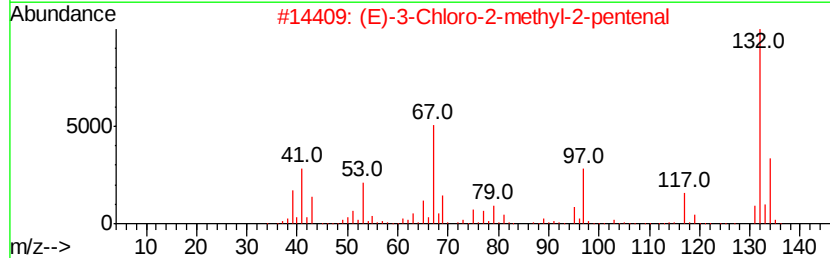
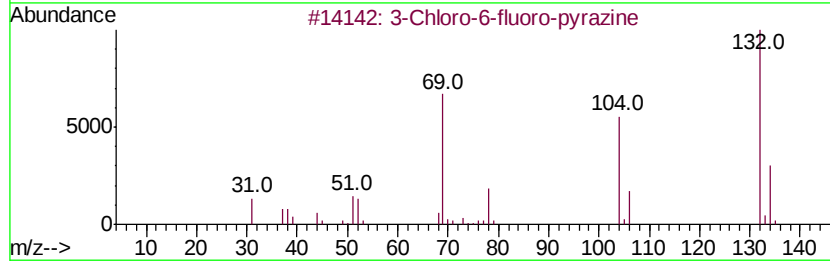
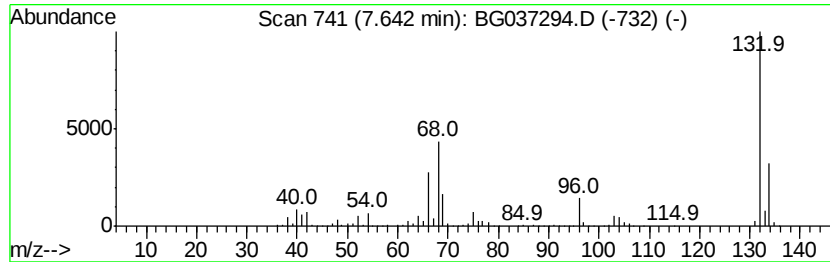
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Peak Number 2 unknown7.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	93.36 ng	1330660	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	3.5	ng	49565	1	8.12	285070	20.0
unknown7.64	7.64	93.4	ng	1330660	1	8.12	285070	20.0